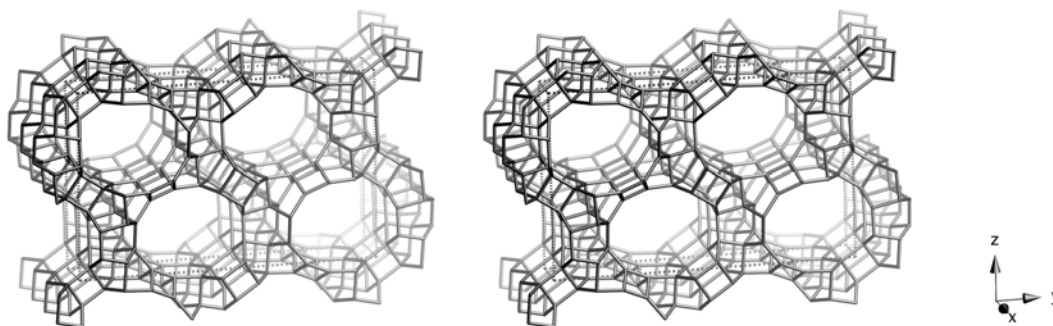


## Framework Type Data



framework viewed along [100]

**Idealized cell data:** orthorhombic, *Cmcm*,  $a = 5.3 \text{ \AA}$ ,  $b = 34.3 \text{ \AA}$ ,  $c = 21.5 \text{ \AA}$

**Coordination sequences and vertex symbols:**

T <sub>1</sub> (8, <i>m</i> ..)	4	11	22	35	57	75	107	135	178	221	269	315	4·6 <sub>2</sub> ·5·6·5·6
T <sub>2</sub> (8, <i>m</i> ..)	4	10	20	31	54	71	106	132	174	212	258	307	4·5·4·5·14 <sub>6</sub> ·*
T <sub>3</sub> (8, <i>m</i> ..)	4	12	22	35	49	80	109	144	178	205	244	321	5·6 <sub>2</sub> ·5·6 <sub>2</sub> ·5 <sub>2</sub> ·6
T <sub>4</sub> (8, <i>m</i> ..)	4	10	19	34	47	76	100	140	172	213	249	299	4·5·4·5·6·14 <sub>6</sub>
T <sub>5</sub> (8, <i>m</i> ..)	4	11	22	38	52	78	103	140	178	220	262	312	4·6 <sub>2</sub> ·5·6·5·6
T <sub>6</sub> (8, <i>m</i> ..)	4	12	19	35	54	79	103	145	167	216	273	315	5·6·5·6·5 <sub>2</sub> ·6
T <sub>7</sub> (8, <i>m</i> ..)	4	12	22	34	50	72	108	147	179	208	248	303	5·6 <sub>2</sub> ·5·6 <sub>2</sub> ·6 <sub>2</sub> ·14 <sub>6</sub>
T <sub>8</sub> (8, <i>m</i> ..)	4	12	23	36	53	75	108	143	178	212	258	314	5·6·5·6·6·6 <sub>2</sub>

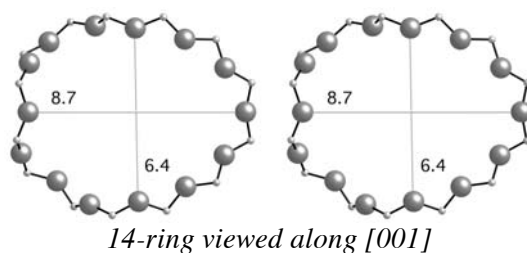
**Secondary building units:** 5-3

**Composite building units:****Materials with this framework type:**

\*SSZ-53<sup>(1)</sup>

## Type Material Data

<b>Crystal chemical data:</b>	[B <sub>1.6</sub> Si <sub>62.4</sub> O <sub>128</sub> ]-SFH monoclinic, C2/c $a = 5.0192\text{\AA}$ , $b = 33.7437\text{\AA}$ , $c = 21.1653\text{\AA}$ , $\beta = 90.485^\circ$ <sup>(1)</sup>
<b>Framework density:</b>	17.9 T/1000Å <sup>3</sup>
<b>Channels:</b>	[001] 14 6.4 x 8.7*

**References:**

- (1) Burton, A., Elomari, S., Chen, C.Y., Medrud, R.C., Chan, I.Y., Bull, L.M., Kibby, C., Harris, T.V., Zones, S.I. and Vittoratos, E.S. *Chem. Eur. Journal*, **9**, 5737-5748 (2003)